

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101619\
 Data File : VN058809.D
 Acq On : 16 Oct 2019 18:02
 Operator : JC/SP
 Sample : K5396-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FA19-JBW7345B

Quant Time: Oct 17 04:45:13 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N101319W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 12 00:00:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	389396	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	616912	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	543781	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	210241	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	289203	53.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.10%	
35) Dibromofluoromethane	7.57	113	205750	54.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.00%	
50) Toluene-d8	10.09	98	775087	50.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.74%	
62) 4-Bromofluorobenzene	12.41	95	246533	42.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	85.48%	
Target Compounds						
16) Acetone	3.80	43	7853	3.480	ug/l	92
27) cis-1,2-Dichloroethene	6.81	96	204945	42.015	ug/l	88
37) Ethyl Acetate	6.92	43	19937	3.161	ug/l #	79
44) Trichloroethene	8.82	130	207063	46.546	ug/l	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN101619\
Data File : VN058809.D
Acq On : 16 Oct 2019 18:02
Operator : JC/SP
Sample : K5396-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FA19-JBW7345B

Quant Time: Oct 17 04:45:13 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N101319W.M
Quant Title : SW846 8260
QLast Update : Sat Oct 12 00:00:30 2019
Response via : Initial Calibration

